

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE072318\
 Data File : BE097030.D
 Acq On : 23 Jul 2018 10:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 23 11:12:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 11:34:20 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	63	0.00
2	1,4-Dioxane	0.573	0.664	-15.9	76	0.00
3	n-Nitrosodimethylamine	0.644	0.675	-4.8	71	0.00
4 S	2-Fluorophenol	0.957	0.851	11.1	64	0.00
5 S	Phenol-d6	1.460	1.191	18.4	57	0.00
6	bis(2-Chloroethyl)ether	1.433	1.347	6.0	62	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	53	0.00
8 S	Nitrobenzene-d5	0.292	0.299	-2.4	59	0.00
9	Naphthalene	3.581	3.535	1.3	55	0.00
10	Hexachlorobutadiene	0.145	0.170	-17.2	66	0.01
11 SURR	2-Methylnaphthalene-d10	0.583	0.550	5.7	53	0.00
12	2-Methylnaphthalene	0.607	0.570	6.1	52	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	57	0.00
14 S	2,4,6-Tribromophenol	0.101	0.083	17.8	57	0.00
15 S	2-Fluorobiphenyl	1.488	1.430	3.9	56	0.00
16	Acenaphthylene	1.827	1.722	5.7	56	0.00
17	Acenaphthene	1.136	1.031	9.2	54	0.00
18	Fluorene	1.275	1.269	0.5	60	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
20	4-Bromophenyl-phenylether	0.189	0.179	5.3	71	0.00
21	Hexachlorobenzene	0.210	0.204	2.9	71	0.00
22	Pentachlorophenol	0.034	0.048	-41.2#	112	0.00
23	Phenanthrene	0.958	0.920	4.0	71	0.00
24	Anthracene	0.834	0.775	7.1	70	0.00
25 SURR	Fluoranthene-d10	3.480	3.785	-8.8	83	0.00
26	Fluoranthene	1.010	1.038	-2.8	77	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	83	0.00
28	Pvrene	1.181	1.072	9.2	78	0.00
29 S	Terphenyl-d14	0.772	0.746	3.4	84	0.00
30	Benzo(a)anthracene	0.952	0.898	5.7	86	0.00
31	Chrysene	1.068	1.021	4.4	82	0.00
32	Bis(2-ethylhexyl)phthalate	0.951	0.633	33.4#	64	0.00
33	Indeno(1,2,3-cd)pyrene	0.778	0.782	-0.5	94	0.00
34 I	Perylene-d12	1.000	1.000	0.0	82	0.00
35	Benzo(b)fluoranthene	1.021	0.869	14.9	75	0.00
36	Benzo(k)fluoranthene	1.214	1.086	10.5	77	0.00
37 C	Benzo(a)pyrene	0.913	0.788	13.7	78	0.00
38	Dibenzo(a,h)anthracene	0.746	0.781	-4.7	97	0.00
39	Benzo(a,h,i)perylene	0.830	0.876	-5.5	96	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE072318\
Data File : BE097030.D
Acq On : 23 Jul 2018 10:16
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 23 11:12:22 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 11:34:20 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			